

s=> fil reg; d stat que l7; fil capl; d que nos l8; fil marpat; d que nos l17
 FILE 'REGISTRY' ENTERED AT 09:54:39 ON 17 DEC 2008
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STRUCTURE FILE UPDATES: 15 DEC 2008 HIGHEST RN 1084993-68-9
 DICTIONARY FILE UPDATES: 15 DEC 2008 HIGHEST RN 1084993-68-9

New CAS Information Use Policies, enter HELP USAGETERMS for details.

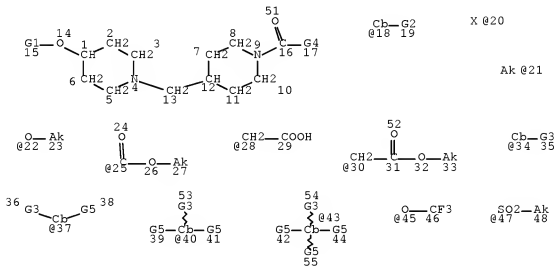
TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

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 conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
 predicted properties as well as tags indicating availability of
 experimental property data in the original document. For information
 on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stdoc/properties.html>

L4 STR



Page 1-A



Page 2-A

VAR G1=56/18
 VAR G2=20/21/22
 VAR G3=OH/COOH/25/28/30
 VAR G4=34/37/40/43

VAR G5=20/CN/OH/21/22/CF3/45/47/49

NODE ATTRIBUTES:

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DEFAULT MLEVEL IS ATOM
MLEVEL IS CLASS AT 18 20 21 23 27 33 34 37 40 43 48 56
DEFAULT ELEVEL IS LIMITED

GRAPH ATTRIBUTES:

RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 56

STEREO ATTRIBUTES: NONE

L7 18 SEA FILE=REGISTRY SSS FUL L4

100.0% PROCESSED 3013 ITERATIONS

18 ANSWERS

SEARCH TIME: 00.00.01

FILE 'CAPLUS' ENTERED AT 09:54:39 ON 17 DEC 2008

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FILE COVERS 1907 - 17 Dec 2008 VOL 149 ISS 25

FILE LAST UPDATED: 16 Dec 2008 (20081216/ED)

Caplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/legal/infopolicy.html>

'OBI' IS DEFAULT SEARCH FIELD FOR 'CAPLUS' FILE

L4 STR

L7 18 SEA FILE=REGISTRY SSS FUL L4

L8 1 SEA FILE=CAPLUS SPE=ON ABB=ON L7

FILE 'MARPAT' ENTERED AT 09:54:39 ON 17 DEC 2008
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FILE CONTENT: 1961-PRESENT VOL 149 ISS 24 (20081212/ED)

MARPAT RECORDS ARE DERIVED FROM INPI DATA FOR 1961-1987

MOST RECENT CITATIONS FOR PATENTS FROM MAJOR ISSUING AGENCIES
 (COVERAGE TO THESE DATES IS NOT COMPLETE):

US	20080280867	13	NOV	2008
DE	102008019744	30	OCT	2008
EP	1990054	12	NOV	2008
JP	2008262895	30	OCT	2008
WO	2008136863	13	NOV	2008
GB	2448808	29	OCT	2008
FR	2915685	07	NOV	2008
RU	2337918	10	NOV	2008
CA	2629177	18	OCT	2008

Expanded G-group definition display now available.

The new MARPAT User Guide is now available at:
<http://www.cas.org/support/stngen/stndoc/marpat.html>.

L4 STR
 L17 1 SEA FILE=MARPAT SSS FUL L4

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FILE 'CAPLUS' ENTERED AT 09:54:44 ON 17 DEC 2008
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FILE 'MARPAT' ENTERED AT 09:54:44 ON 17 DEC 2008
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 COPYRIGHT (C) 2008 American Chemical Society (ACS)
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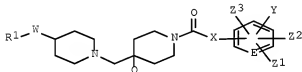
=> d ibib abs hitstr l18; fil hom

L18 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2008 ACS on STN DUPLICATE 1
 ACCESSION NUMBER: 2004:817881 CAPLUS Full-text
 DOCUMENT NUMBER: 141:332055
 TITLE: Preparation of piperidine derivatives for the
 treatment of chemokine or H1 mediated disease state
 INVENTOR(S): Luckhurst, Christopher; Perry,
 Matthew; Springthorpe, Brian
 PATENT ASSIGNEE(S): Astrazeneca AB, Swed.

SOURCE: PCT Int. Appl., 44 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004085423	A1	20041007	WO 2004-SE450	20040323
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EP 1611124	A1	20060104	EP 2004-722752	20040323
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK				
JP 2006521374	T	20060921	JP 2006-507982	20040323
US 20060281726	A1	20061214	US 2005-549868	20050921
PRIORITY APPLN. INFO.:			SE 2003-850	A 20030325
			WO 2004-SE450	W 20040323
OTHER SOURCE(S):		MARPAT 141:332055		

GI

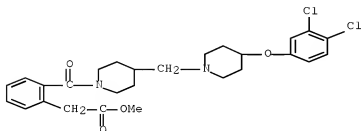


I

- AB The title compds. [I; E = CH, N; Q = H, OH; W = CH₂, O, NR₂; X = a bond, CH₂, CH₂O; Y = OH, SO₃H, etc.; Z₁-Z₃ = H, halo, CN, NO₂, etc.; R₁ = (un)substituted Ph; R₂ = H, alkyl], useful in the treatment of a chemokine (such as CCR3) or H1 mediated disease state, were prepared. Thus, reacting 4-[[4-(3,4-dichlorophenoxy)piperidin-1-yl]methyl]piperidine with phthalic anhydride followed by treatment of the reaction mixture with AcOH afforded 2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoic acid which showed pK_i of 6.5 in human H1 receptor binding assay. The pharmaceutical composition comprising the compound I is claimed.
- IT 779729-78-7P, Methyl 2-[2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]phenyl]acetate
 779729-81-2P, Methyl 4-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoate
 RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
 (preparation of piperidine derivs. for the treatment of chemokine or H1 mediated disease state)

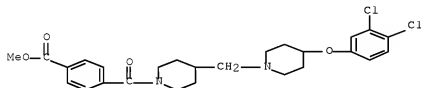
RN 770729-78-7 CAPLUS

CN Benzenecetic acid, 2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-, methyl ester (CA INDEX NAME)



RN 770729-81-2 CAPLUS

CN Benzoic acid, 4-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-, methyl ester (CA INDEX NAME)



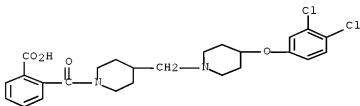
IT 770729-73-2P, 2-[[4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoic acid
 770729-74-3P, 2-[[4-[[4-(2,4-Dichloro-3-methylphenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoic acid
 770729-75-4P, 2-[[4-[[4-(3,4-Dichloro-2-methylphenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoic acid
 770729-76-5P, 2-[[4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-3,6-difluorobenzoic acid
 770729-79-8P, Methyl 3-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoate 770729-80-1P
 , Methyl 2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-4-methoxybenzoate 770729-83-4P, Methyl
 4-chloro-2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoate 770729-84-5P,
 2-[2-[[4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]phenyl]acetic acid 770729-85-6P,
 3-[[4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoic acid 770729-87-8P,
 2-[[4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-4-methoxybenzoic acid 770729-88-9P,
 4-[[4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoic acid 770729-89-0P,
 2-[[4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-4-methylbenzoic acid 770729-90-3P,
 4-Chloro-2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]benzoic acid sodium salt 770729-91-4P,
 4-[[4-(3,4-Dichlorophenoxy)-1-piperidinyl]methyl]-1-[4-hydroxy-3-(methylsulfonyl)benzoyl]piperidine

RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(preparation of piperidine derivs. for the treatment of chemokine or H1 mediated disease state)

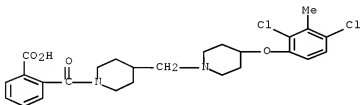
RN 770729-73-2 CAPLUS

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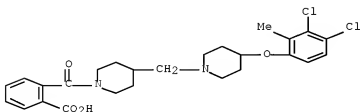
RN 770729-74-3 CAPLUS

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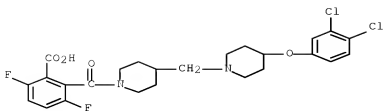
RN 770729-75-4 CAPLUS

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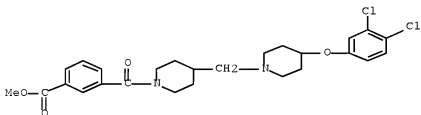
RN 770729-76-5 CAPLUS

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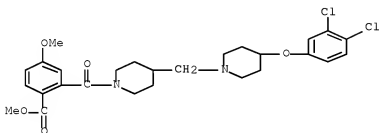
RN 770729-79-8 CAPLUS

CN Benzoic acid, 3-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-, methyl ester (CA INDEX NAME)



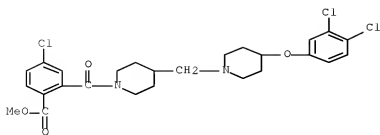
RN 770729-80-1 CAPLUS

CN Benzoic acid, 2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-4-methoxy-, methyl ester (CA INDEX NAME)



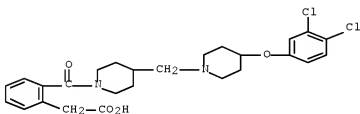
RN 770729-83-4 CAPLUS

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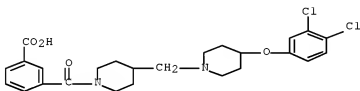
RN 770729-84-5 CAPLUS

CN Benzoic acid, 2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]- (CA INDEX NAME)



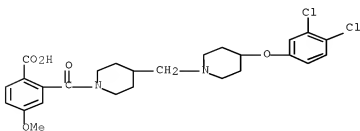
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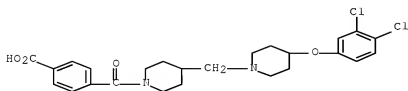
RN 770729-87-8 CAPLUS

CN Benzoic acid, 2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-4-methoxy- (CA INDEX NAME)



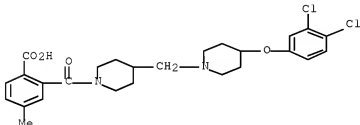
RN 770729-88-9 CAPLUS

CN Benzoic acid, 4-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]- (CA INDEX NAME)



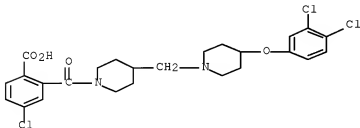
RN 770729-89-0 CAPLUS

CN Benzoic acid, 2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-4-methyl- (CA INDEX NAME)



RN 770729-90-3 CAPLUS

CN Benzoic acid, 4-chloro-2-[[4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl]carbonyl]-, sodium salt (1:1) (CA INDEX NAME)



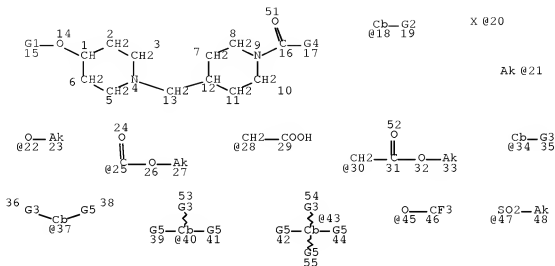
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RN 770729-91-4 CAPLUS

CN Methanone, [4-[[4-(3,4-dichlorophenoxy)-1-piperidinyl]methyl]-1-piperidinyl][4-hydroxy-3-(methylsulfonyl)phenyl]- (CA INDEX NAME)

SEARCH HISTORY

=> d stat que l7; d his nofile
L4 STR



Page 1-A



Ph 056

Page 2-A

VAR G1=56/18
VAR G2=20/21/22
VAR G3=OH/COOH/25/28/30
VAR G4=34/37/40/43
VAR G5=20/CN/OH/21/22/CF3/45/47/49
NODE ATTRIBUTES:
CONNECT IS E1 RC AT 21
CONNECT IS E1 RC AT 23
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CONNECT IS E4 RC AT 40
CONNECT IS E5 RC AT 43
CONNECT IS E1 RC AT 48
DEFAULT MLEVEL IS ATOM
MLEVEL IS CLASS AT 18 20 21 23 27 33 34 37 40 43 48 56
DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:
RING(S) ARE ISOLATED OR EMBEDDED
NUMBER OF NODES IS 56

STEREO ATTRIBUTES: NONE
L7 18 SEA FILE=REGISTRY SSS FUL L4

100.0% PROCESSED 3013 ITERATIONS
SEARCH TIME: 00.00.01

18 ANSWERS

(FILE 'HOME' ENTERED AT 09:32:47 ON 17 DEC 2008)

FILE 'CAPLUS' ENTERED AT 09:33:03 ON 17 DEC 2008
E US2006-549868/APPS

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D SCAN
SEL RN

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L7 18 SEA SSS FUL L4
SAVE TEMP L7 CHA868FULL/A

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L8 1 SEA SPE=ON ABB=ON L7

L9 20 SEA SPE=ON ABB=ON LUCKHURST C?/AU
 L10 1565 SEA SPE=ON ABB=ON PERRY M?/AU
 L11 48 SEA SPE=ON ABB=ON SPRINGTHORPE B?/AU
 L12 1 SEA SPE=ON ABB=ON (L9 OR L10 OR L11) AND L8
 L13 0 SEA SPE=ON ABB=ON L1 AND L8
 D SCAN TI L12
 D SCAN TI L1

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 ANALYZE L7 1- LC : 3 TERMS
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FILE 'STNGUIDE' ENTERED AT 09:49:48 ON 17 DEC 2008

L15 FILE 'MARPAT' ENTERED AT 09:50:01 ON 17 DEC 2008
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 L16 12645 SEA SSS FUL L4 EXTEND
 L17 1 SEA SSS FUL L4
 SAVE TEMP L17 CHA868MARPA
 D SCAN

FILE 'STNGUIDE' ENTERED AT 09:53:53 ON 17 DEC 2008

FILE 'REGISTRY' ENTERED AT 09:54:39 ON 17 DEC 2008
 D STAT QUE L7

FILE 'CAPLUS' ENTERED AT 09:54:39 ON 17 DEC 2008
 D QUE NOS L8

FILE 'MARPAT' ENTERED AT 09:54:39 ON 17 DEC 2008
 D QUE NOS L17

L18 FILE 'CAPLUS, MARPAT' ENTERED AT 09:54:44 ON 17 DEC 2008
 1 DUP REM L12 L17 (1 DUPLICATE REMOVED)
 ANSWER '1' FROM FILE CAPLUS
 D IBIB ABS HITSTR L18

FILE 'HOME' ENTERED AT 09:54:54 ON 17 DEC 2008
 D STAT QUE L7

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